

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101019\
 Data File : BF117367.D
 Acq On : 10 Oct 2019 18:29
 Operator : JU
 Sample : K5297-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Quant Time: Oct 11 06:46:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:26:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	45710	20.00	ng	-0.01
21) Naphthalene-d8	8.07	136	181719	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	94464	20.00	ng	-0.01
64) Phenanthrene-d10	11.32	188	149077	20.00	ng	-0.01
76) Chrysene-d12	13.96	240	85974	20.00	ng	-0.01
87) Perylene-d12	15.42	264	98624	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	284864	97.30	ng	0.00
7) Phenol-d6	6.44	99	383782	101.43	ng	-0.01
23) Nitrobenzene-d5	7.36	82	232031	67.09	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	76845	97.33	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	403998	66.87	ng	-0.01
79) Terphenyl-d14	12.90	244	315586	68.61	ng	-0.01
Target Compounds						
10) Phenol	6.45	94	15176	3.638	ng	93
31) Naphthalene	8.10	128	23082	2.641	ng	# 92
37) 2-Methylnaphthalene	8.79	142	30165	5.126	ng	97
38) 1-Methylnaphthalene	8.89	142	19986	3.626	ng	90
50) Dimethylphthalate	9.55	163	45641	6.824	ng	99
52) Acenaphthene	9.86	154	25498	4.910	ng	96
55) Dibenzofuran	10.04	168	23658	3.088	ng	96
58) Fluorene	10.38	166	25900	4.196	ng	# 96
71) Phenanthrene	11.34	178	107804	12.731	ng	98
75) Fluoranthene	12.53	202	66995	7.700	ng	96
78) Pyrene	12.76	202	44271	6.137	ng	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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